



Full wwPDB NMR Structure Validation Report ⓘ

May 28, 2026 – 09:19 AM EDT

PDB ID : 9K97 / pdb_00009k97
BMRB ID : 52617
Title : Solution structure of DRB2 dsRBD2
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Deposited on : 2024-10-25

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with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
wwPDB-RCI : v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV : Wang et al. (2010)
wwPDB-ShiftChecker : v1.2
BMRB Restraints Analysis : v1.2
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

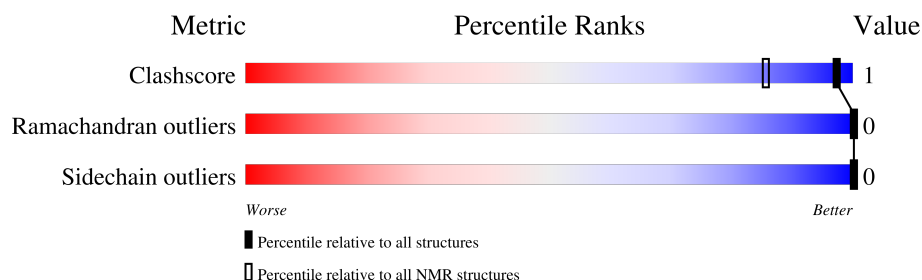
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR


The overall completeness of chemical shifts assignment is 32%.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	NMR archive (#Entries)
Clashscore	229148	14424
Ramachandran outliers	224038	12848
Sidechain outliers	223484	12823

The table below summarises the geometric issues observed across the polymeric chains and their fit to the experimental data. The red, orange, yellow and green segments indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A cyan segment indicates the fraction of residues that are not part of the well-defined cores, and a grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$

Mol	Chain	Length	Quality of chain
1	A	100	

2 Ensemble composition and analysis

This entry contains 10 models. Model 7 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative, based on the following criterion: *lowest energy*.

The following residues are included in the computation of the global validation metrics.

Well-defined (core) protein residues			
Well-defined core	Residue range (total)	Backbone RMSD (Å)	Medoid model
1	A:87-A:156, A:173-A:186 (84)	0.88	7

Ill-defined regions of proteins are excluded from the global statistics.

Ligands and non-protein polymers are included in the analysis.

The models can be grouped into 2 clusters and 1 single-model cluster was found.

Cluster number	Models
1	1, 2, 4, 5, 7, 9, 10
2	3, 8
Single-model clusters	6

3 Entry composition [i](#)

There is only 1 type of molecule in this entry. The entry contains 1565 atoms, of which 789 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Double-stranded RNA-binding protein 2.

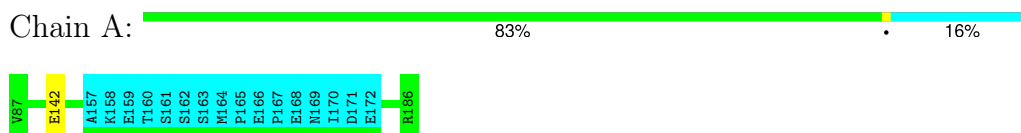
Mol	Chain	Residues	Atoms						Trace
1	A	100	Total	C	H	N	O	S	0
			1565	490	789	136	148	2	

4 Residue-property plots

4.1 Average score per residue in the NMR ensemble

These plots are provided for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic is the same as shown in the summary in section 1 of this report. The second graphic shows the sequence where residues are colour-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. Stretches of 2 or more consecutive residues without any outliers are shown as green connectors. Residues which are classified as ill-defined in the NMR ensemble, are shown in cyan with an underline colour-coded according to the previous scheme. Residues which were present in the experimental sample, but not modelled in the final structure are shown in grey.

- Molecule 1: Double-stranded RNA-binding protein 2

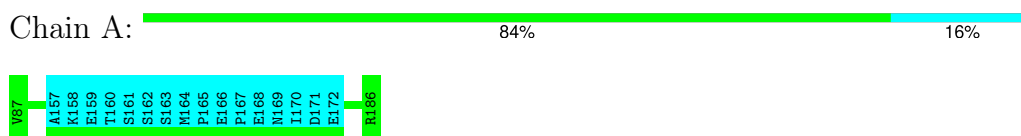


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

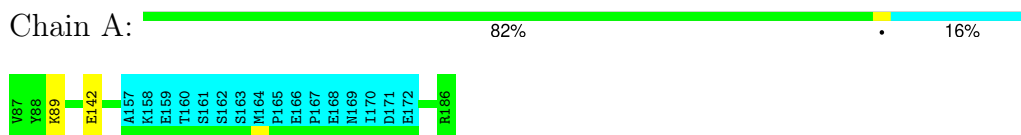
4.2.1 Score per residue for model 1

- Molecule 1: Double-stranded RNA-binding protein 2



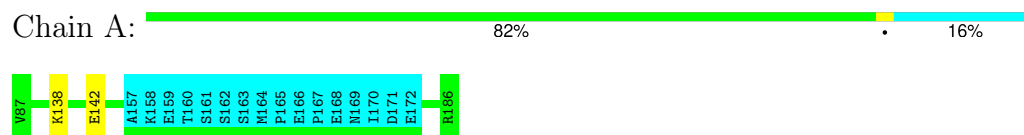
4.2.2 Score per residue for model 2

- Molecule 1: Double-stranded RNA-binding protein 2



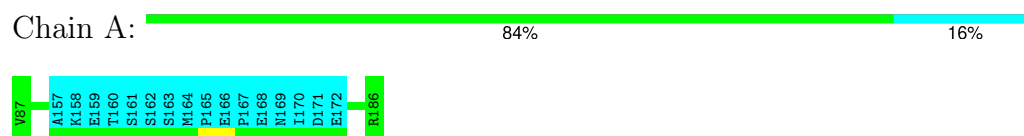
4.2.3 Score per residue for model 3

- Molecule 1: Double-stranded RNA-binding protein 2



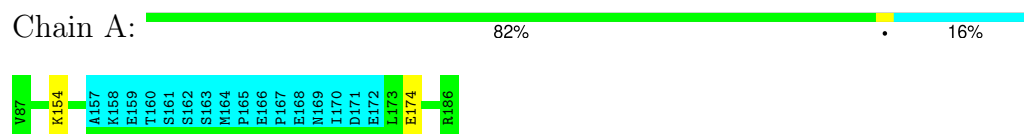
4.2.4 Score per residue for model 4

- Molecule 1: Double-stranded RNA-binding protein 2



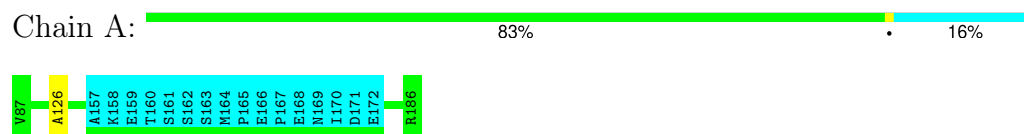
4.2.5 Score per residue for model 5

- Molecule 1: Double-stranded RNA-binding protein 2



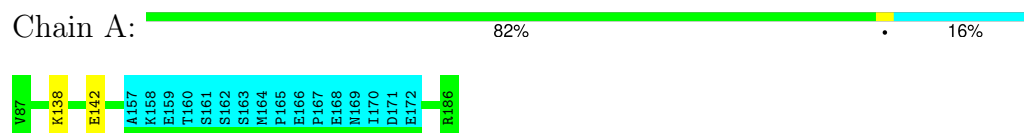
4.2.6 Score per residue for model 6

- Molecule 1: Double-stranded RNA-binding protein 2



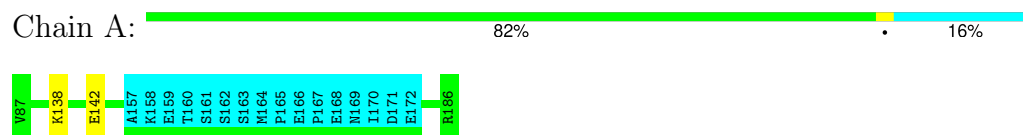
4.2.7 Score per residue for model 7 (medoid)

- Molecule 1: Double-stranded RNA-binding protein 2



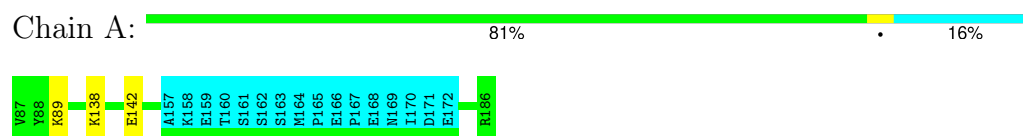
4.2.8 Score per residue for model 8

- Molecule 1: Double-stranded RNA-binding protein 2



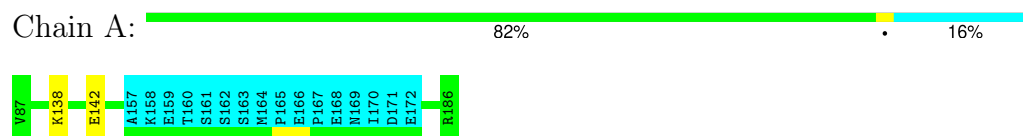
4.2.9 Score per residue for model 9

- Molecule 1: Double-stranded RNA-binding protein 2



4.2.10 Score per residue for model 10

- Molecule 1: Double-stranded RNA-binding protein 2



5 Refinement protocol and experimental data overview

The models were refined using the following method: *torsion angle dynamics*.

Of the 5000 calculated structures, 10 were deposited, based on the following criterion: *structures with the lowest energy*.

The following table shows the software used for structure solution, optimisation and refinement.

Software name	Classification	Version
CS-RDC-NOE Rosetta	structure calculation	
CS-RDC-NOE Rosetta	refinement	

The following table shows chemical shift validation statistics as aggregates over all chemical shift files. Detailed validation can be found in section 7 of this report.

Chemical shift file(s)	working_cs.cif
Number of chemical shift lists	1
Total number of shifts	474
Number of shifts mapped to atoms	434
Number of unparsed shifts	0
Number of shifts with mapping errors	40
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	32%

6 Model quality [i](#)

6.1 Standard geometry [i](#)

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the (average) root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#Z>5	RMSZ	#Z>5
1	A	0.47±0.00	0±0/667 (0.0± 0.0%)	0.65±0.01	0±0/901 (0.0± 0.0%)
All	All	0.47	0/6670 (0.0%)	0.65	1/9010 (0.0%)

There are no bond-length outliers.

All unique angle outliers are listed below.

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	126	ALA	CB-CA-C	-5.48	110.24	116.54	6	1

There are no chirality outliers.

There are no planarity outliers.

6.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in each chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes averaged over the ensemble.

Mol	Chain	Non-H	H(model)	H(added)	Clashes
1	A	655	681	679	1±1
All	All	6550	6810	6790	8

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 1.

All unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:154:LYS:NZ	1:A:174:GLU:OE1	0.49	2.43	5	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:89:LYS:NZ	1:A:142:GLU:OE1	0.48	2.42	2	2
1:A:138:LYS:NZ	1:A:142:GLU:OE2	0.46	2.48	10	5

6.3 Torsion angles [i](#)

6.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all NMR entries. The Analysed column shows the number of residues for which the backbone conformation was analysed and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	82/100 (82%)	81±1 (99±1%)	1±1 (1±1%)	0±0 (0±0%)	100	100
All	All	820/1000 (82%)	814 (99%)	6 (1%)	0 (0%)	100	100

There are no Ramachandran outliers.

6.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all NMR entries. The Analysed column shows the number of residues for which the sidechain conformation was analysed and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	67/82 (82%)	67±0 (100±0%)	0±0 (0±0%)	100	100
All	All	670/820 (82%)	670 (100%)	0 (0%)	100	100

There are no protein residues with a non-rotameric sidechain to report.

6.3.3 RNA [i](#)

There are no RNA molecules in this entry.

6.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.6 Ligand geometry [i](#)

There are no ligands in this entry.

6.7 Other polymers [i](#)

There are no such molecules in this entry.

6.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

7 Chemical shift validation

The completeness of assignment taking into account all chemical shift lists is 32% for the well-defined parts and 32% for the entire structure.

7.1 Chemical shift list 1

File name: working_cs.cif

Chemical shift list name: *default*

7.1.1 Bookkeeping

The following table shows the results of parsing the chemical shift list and reports the number of nuclei with statistically unusual chemical shifts.

Total number of shifts	474
Number of shifts mapped to atoms	434
Number of unparsed shifts	0
Number of shifts with mapping errors	40
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	0

The following assigned chemical shifts were not mapped to the molecules present in the coordinate file.

- No matching atom found in the structure. All 40 occurrences are reported below.

List ID	Chain	Res	Type	Atom	Shift Data		
					Value	Uncertainty	Ambiguity
1	A	85	THR	C	175.0064223	.	1
1	A	85	THR	CA	62.57291323	.	1
1	A	85	THR	CB	68.88254714	.	1
1	A	86	GLY	C	174.520548	0.02443703815	1
1	A	86	GLY	CA	45.68234614	0.03013691385	1
1	A	86	GLY	H	8.449894731	0.001101206864	1
1	A	86	GLY	N	112.6104357	0.3998618864	1
1	A	87	VAL	H	7.63194556	0.001831199846	1
1	A	189	GLU	C	178.0274722	.	1
1	A	189	GLU	CA	56.76353455	.	1
1	A	189	GLU	CB	29.26501362	.	1
1	A	190	ASN	C	174.9624098	0.01089714312	1
1	A	190	ASN	CA	52.43204636	0.330543899	1
1	A	190	ASN	CB	38.29426341	0.08126574848	1

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List ID	Chain	Res	Type	Atom	Shift Data		
					Value	Uncertainty	Ambiguity
1	A	190	ASN	H	8.1511781	0.003222972115	1
1	A	190	ASN	N	119.663192	0.3579786335	1
1	A	191	ILE	C	176.2778538	0.3633745715	1
1	A	191	ILE	CA	60.94215442	0.05729919968	1
1	A	191	ILE	CB	37.62565243	0.2421469117	1
1	A	191	ILE	H	7.95158245	0.001203355583	1
1	A	191	ILE	N	122.4638897	0.3511568964	1
1	A	193	THR	H	7.896785986	0.001522046731	1
1	A	193	THR	N	114.2726699	0.3428360899	1
1	A	195	SER	C	174.5422161	0.002621855889	1
1	A	195	SER	CA	58.25539253	0.2685598871	1
1	A	195	SER	CB	63.37777544	0.15368116	1
1	A	195	SER	H	8.080172704	0.001150149141	1
1	A	195	SER	N	117.2356775	0.3926799424	1
1	A	197	LEU	C	177.1733283	0.008373092571	1
1	A	197	LEU	CA	54.94371086	0.01121952147	1
1	A	197	LEU	CB	40.90556494	0.01496494195	1
1	A	197	LEU	H	8.083867149	0.002239043355	1
1	A	197	LEU	N	125.023748	0.3041954199	1
1	A	198	GLU	H	8.051984628	0.000889371177	1
1	A	198	GLU	N	121.9907159	0.3203393562	1
1	A	199	HIS	C	179.0642109	0.003895081913	1
1	A	199	HIS	CA	56.87969865	.	1
1	A	199	HIS	CB	29.66684177	.	1
1	A	199	HIS	H	7.873354169	0.003792075446	1
1	A	199	HIS	N	126.9744263	0.2971056295	1

7.1.2 Chemical shift referencing [i](#)

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction \pm precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	99	0.42 ± 0.07	None needed (< 0.5 ppm)
$^{13}\text{C}_\beta$	93	1.18 ± 0.11	Should be checked
$^{13}\text{C}'$	98	0.07 ± 0.06	None needed (< 0.5 ppm)
^{15}N	92	-1.61 ± 0.15	Should be applied

7.1.3 Completeness of resonance assignments [i](#)

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 32%, i.e. 371 atoms were assigned a chemical

shift out of a possible 1172. 0 out of 14 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹H	¹³C	¹⁵N
Backbone	298/418 (71%)	72/170 (42%)	154/168 (92%)	72/80 (90%)
Sidechain	73/679 (11%)	0/444 (0%)	73/204 (36%)	0/31 (0%)
Aromatic	0/75 (0%)	0/37 (0%)	0/37 (0%)	0/1 (0%)
Overall	371/1172 (32%)	72/651 (11%)	227/409 (56%)	72/112 (64%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 32%, i.e. 435 atoms were assigned a chemical shift out of a possible 1360. 0 out of 14 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹H	¹³C	¹⁵N
Backbone	349/494 (71%)	84/200 (42%)	181/200 (90%)	84/94 (89%)
Sidechain	86/791 (11%)	0/515 (0%)	86/243 (35%)	0/33 (0%)
Aromatic	0/75 (0%)	0/37 (0%)	0/37 (0%)	0/1 (0%)
Overall	435/1360 (32%)	84/752 (11%)	267/480 (56%)	84/128 (66%)

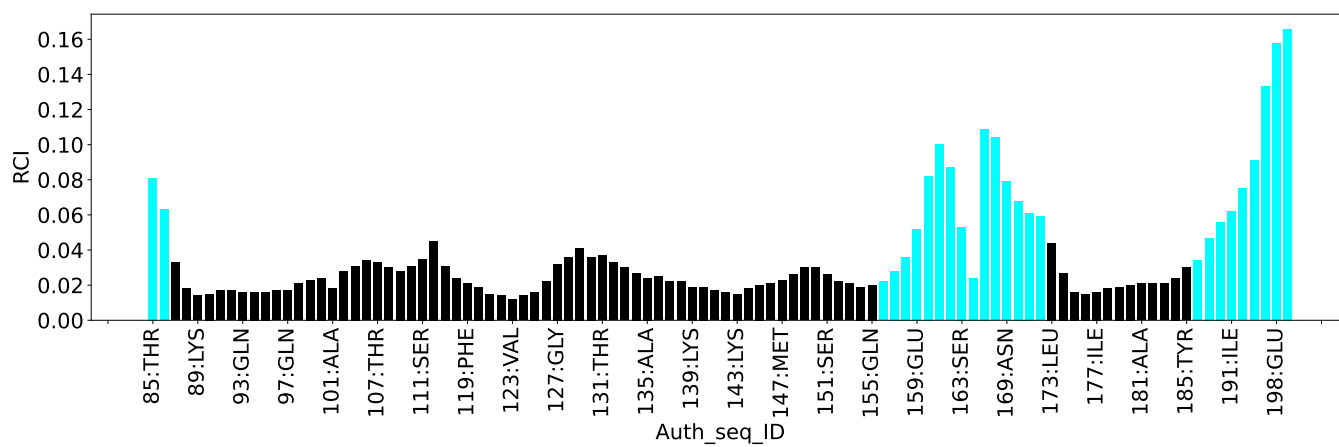
7.1.4 Statistically unusual chemical shifts [i](#)

There are no statistically unusual chemical shifts.

7.1.5 Random Coil Index (RCI) plots [i](#)

The image below reports *random coil index* values for the protein chains in the structure. The height of each bar gives a probability of a given residue to be disordered, as predicted from the available chemical shifts and the amino acid sequence. A value above 0.2 is an indication of significant predicted disorder. The colour of the bar shows whether the residue is in the well-defined core (black) or in the ill-defined residue ranges (cyan), as described in section 2 on ensemble composition. If well-defined core and ill-defined regions are not identified then it is shown as gray bars.

Random coil index (RCI) for chain A:



8 NMR restraints analysis

8.1 Conformationally restricting restraints

The following table provides the summary of experimentally observed NMR restraints in different categories. Restraints are classified into different categories based on the sequence separation of the atoms involved.

Description	Value
Total distance restraints	35
Intra-residue ($ i-j =0$)	0
Sequential ($ i-j =1$)	0
Medium range ($ i-j >1$ and $ i-j <5$)	0
Long range ($ i-j \geq 5$)	35
Inter-chain	0
Hydrogen bond restraints	0
Disulfide bond restraints	0
Total dihedral-angle restraints	0
Number of unmapped restraints	0
Number of restraints per residue	0.3
Number of long range restraints per residue ¹	0.3

¹Long range hydrogen bonds and disulfide bonds are counted as long range restraints while calculating the number of long range restraints per residue

8.2 Residual restraint violations

This section provides the overview of the restraint violations analysis. The violations are binned as small, medium and large violations based on its absolute value. Average number of violations per model is calculated by dividing the total number of violations in each bin by the size of the ensemble.

8.2.1 Average number of distance violations per model

Distance violations less than 0.1 Å are not included in the calculation.

Bins (Å)	Average number of violations per model	Max (Å)
0.1-0.2 (Small)	0.2	0.13
0.2-0.5 (Medium)	0.4	0.5
>0.5 (Large)	0.2	0.84

8.2.2 Average number of dihedral-angle violations per model [i](#)

Dihedral-angle violations less than 1° are not included in the calculation. There are no dihedral-angle violations

9 Distance violation analysis

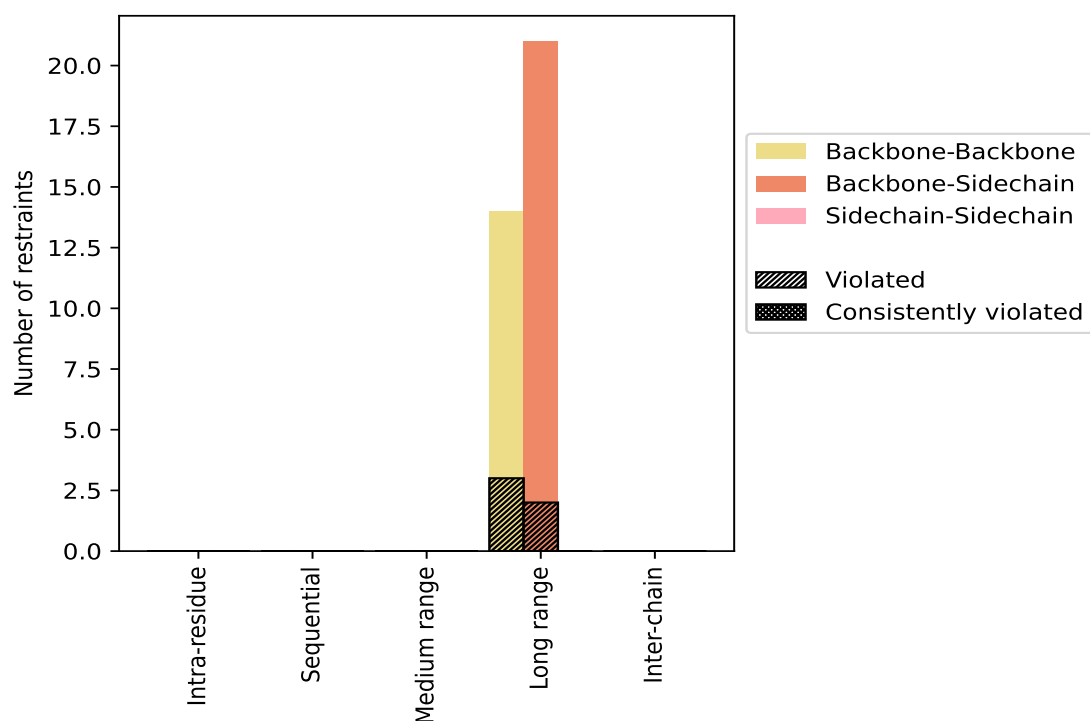
9.1 Summary of distance violations

The following table shows the summary of distance violations in different restraint categories based on the sequence separation of the atoms involved. Each category is further sub-divided into three sub-categories based on the atoms involved. Violations less than 0.1 Å are not included in the statistics.

Restrains type	Count	% ¹	Violated ³			Consistently Violated ⁴		
			Count	% ²	% ¹	Count	% ²	% ¹
Intra-residue (i-j =0)	0	0.0	0	0.0	0.0	0	0.0	0.0
Backbone-Backbone	0	0.0	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Sidechain-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Sequential (i-j =1)	0	0.0	0	0.0	0.0	0	0.0	0.0
Backbone-Backbone	0	0.0	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Sidechain-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Medium range (i-j >1 & i-j <5)	0	0.0	0	0.0	0.0	0	0.0	0.0
Backbone-Backbone	0	0.0	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Sidechain-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Long range (i-j ≥5)	35	100.0	5	14.3	14.3	0	0.0	0.0
Backbone-Backbone	14	40.0	3	21.4	8.6	0	0.0	0.0
Backbone-Sidechain	21	60.0	2	9.5	5.7	0	0.0	0.0
Sidechain-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Inter-chain	0	0.0	0	0.0	0.0	0	0.0	0.0
Backbone-Backbone	0	0.0	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Sidechain-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Hydrogen bond	0	0.0	0	0.0	0.0	0	0.0	0.0
Disulfide bond	0	0.0	0	0.0	0.0	0	0.0	0.0
Total	35	100.0	5	14.3	14.3	0	0.0	0.0
Backbone-Backbone	14	40.0	3	21.4	8.6	0	0.0	0.0
Backbone-Sidechain	21	60.0	2	9.5	5.7	0	0.0	0.0
Sidechain-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0

¹ percentage calculated with respect to the total number of distance restraints, ² percentage calculated with respect to the number of restraints in a particular restraint category, ³ violated in at least one model, ⁴ violated in all the models

9.1.1 Bar chart : Distribution of distance restraints and violations [i](#)



Violated and consistently violated restraints are shown using different hatch patterns in their respective categories. The hydrogen bonds and disulfied bonds are counted in their appropriate category on the x-axis

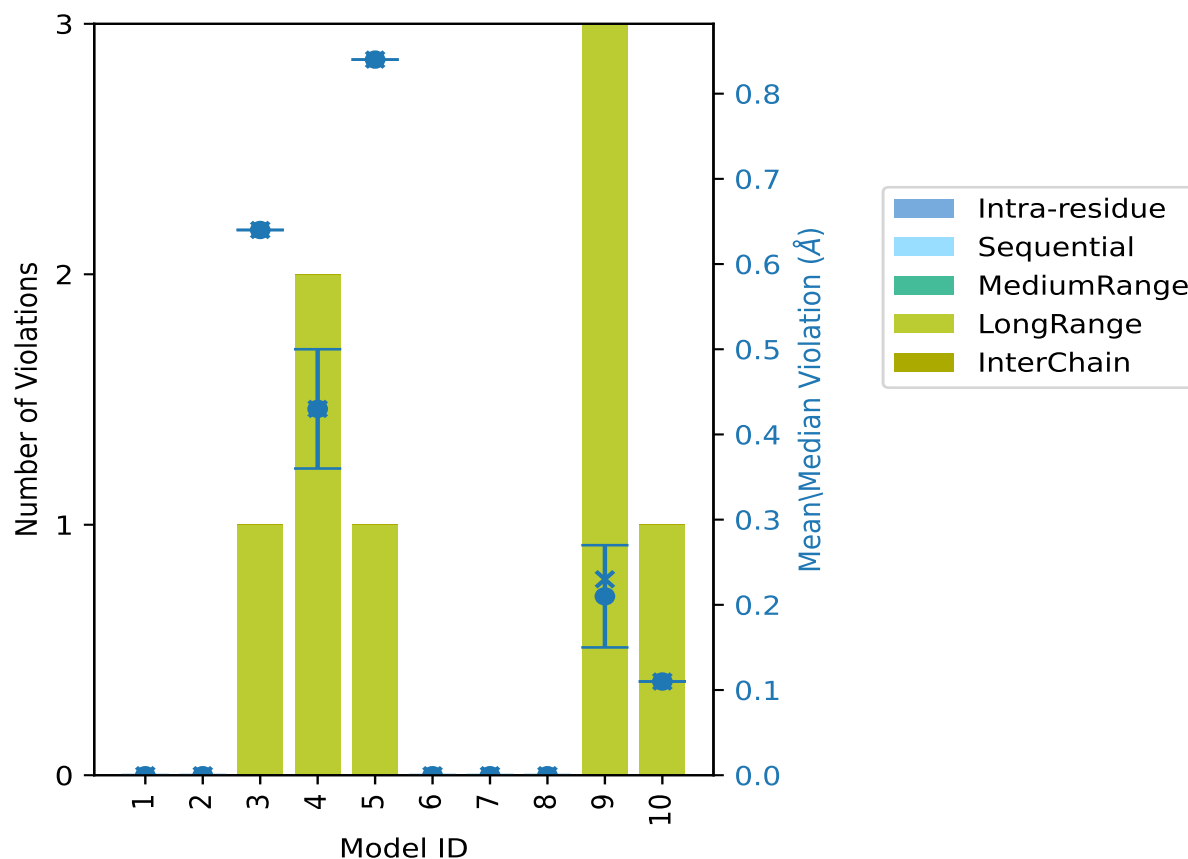
9.2 Distance violation statistics for each model [i](#)

The following table provides the distance violation statistics for each model in the ensemble. Violations less than 0.1 Å are not included in the statistics.

Model ID	Number of violations						Mean (Å)	Max (Å)	SD ⁶ (Å)	Median (Å)
	IR ¹	SQ ²	MR ³	LR ⁴	IC ⁵	Total				
1	0	0	0	0	0	0	0.0	0.0	0.0	0.0
2	0	0	0	0	0	0	0.0	0.0	0.0	0.0
3	0	0	0	1	0	1	0.64	0.64	0.0	0.64
4	0	0	0	2	0	2	0.43	0.5	0.07	0.43
5	0	0	0	1	0	1	0.84	0.84	0.0	0.84
6	0	0	0	0	0	0	0.0	0.0	0.0	0.0
7	0	0	0	0	0	0	0.0	0.0	0.0	0.0
8	0	0	0	0	0	0	0.0	0.0	0.0	0.0
9	0	0	0	3	0	3	0.21	0.27	0.06	0.23
10	0	0	0	1	0	1	0.11	0.11	0.0	0.11

¹Intra-residue restraints, ²Sequential restraints, ³Medium range restraints, ⁴Long range restraints, ⁵Inter-chain restraints, ⁶Standard deviation

9.2.1 Bar graph : Distance Violation statistics for each model [i](#)



The mean(dot),median(x) and the standard deviation are shown in blue with respect to the y axis on the right

9.3 Distance violation statistics for the ensemble [i](#)

Violation analysis may find that some restraints are violated in few models and some are violated in most of models. The following table provides this information as number of violated restraints for a given fraction of the ensemble. In total, 30(IR:0, SQ:0, MR:0, LR:30, IC:0) restraints are not violated in the ensemble.

Number of violated restraints						Fraction of the ensemble	
IR ¹	SQ ²	MR ³	LR ⁴	IC ⁵	Total	Count ⁶	%
0	0	0	2	0	2	1	10.0
0	0	0	3	0	3	2	20.0
0	0	0	0	0	0	3	30.0

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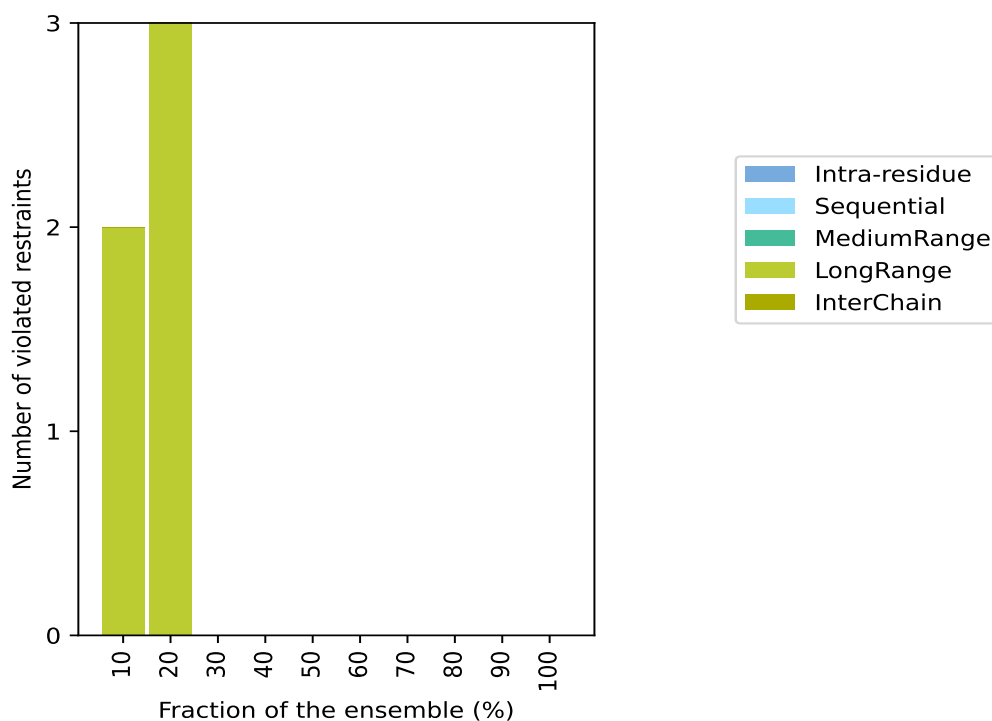
Continued from previous page...

Number of violated restraints						Fraction of the ensemble	
IR ¹	SQ ²	MR ³	LR ⁴	IC ⁵	Total	Count ⁶	%
0	0	0	0	0	0	4	40.0
0	0	0	0	0	0	5	50.0
0	0	0	0	0	0	6	60.0
0	0	0	0	0	0	7	70.0
0	0	0	0	0	0	8	80.0
0	0	0	0	0	0	9	90.0
0	0	0	0	0	0	10	100.0

¹Intra-residue restraints, ²Sequential restraints, ³Medium range restraints, ⁴Long range restraints,

⁵Inter-chain restraints, ⁶ Number of models with violations

9.3.1 Bar graph : Distance violation statistics for the ensemble ⓘ

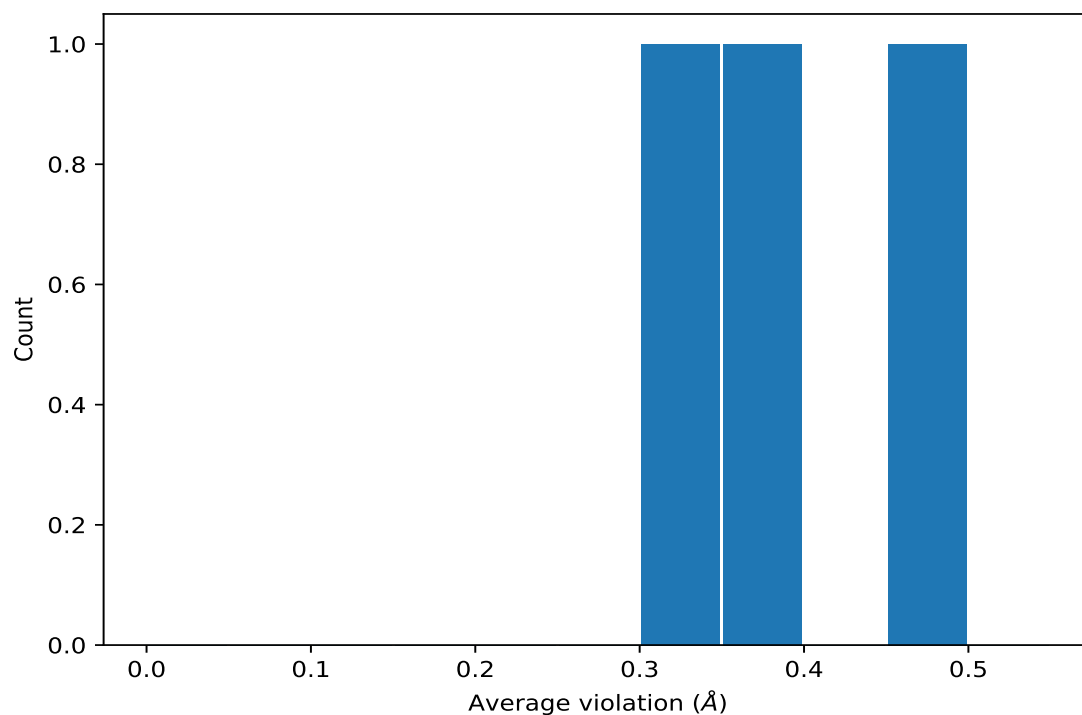


9.4 Most violated distance restraints in the ensemble ⓘ

9.4.1 Histogram : Distribution of mean distance violations ⓘ

The following histogram shows the distribution of the average value of the violation. The average is calculated for each restraint that is violated in more than one model over all the violated models

in the ensemble



9.4.2 Table: Most violated distance restraints [i](#)

The following table provides the mean and the standard deviation of the violation for each restraint sorted by number of violated models and the mean value. The Key (restraint list ID, restraint ID) is the unique identifier for a given restraint. Rows with same key represent combinatorial or ambiguous restraints and are counted as a single restraint.

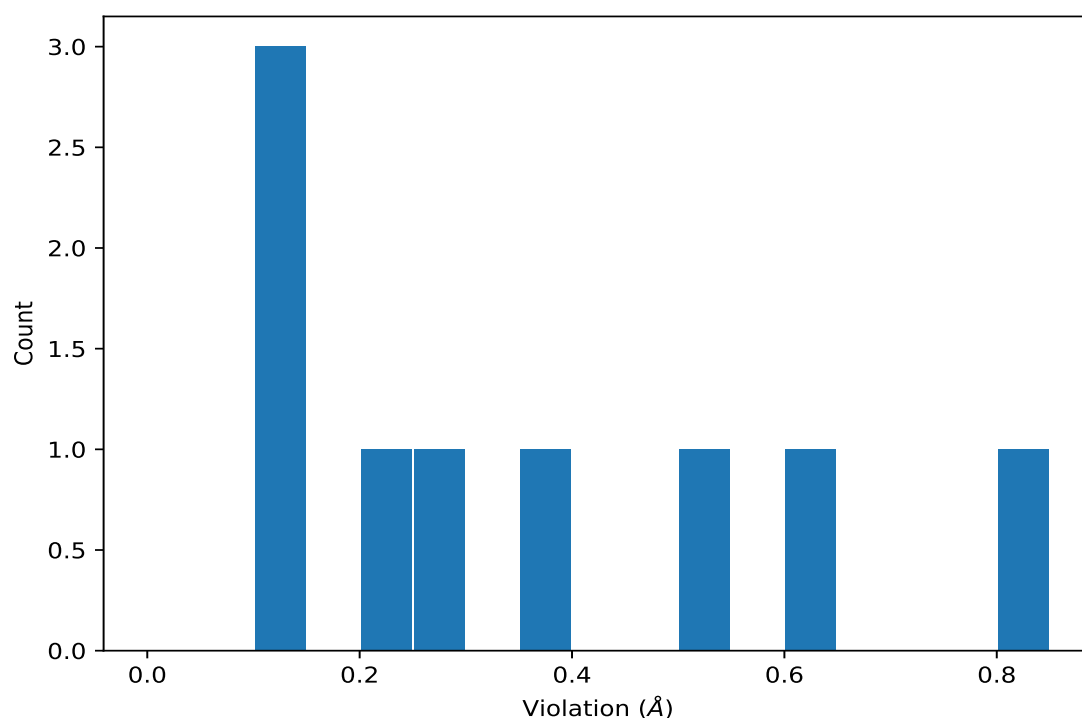
Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(1,6)	1:181:A:ALA:H	1:157:A:ALA:HA	2	0.48	0.36	0.48
(1,29)	1:122:A:THR:H	1:145:A:ALA:HA	2	0.36	0.13	0.36
(1,33)	1:133:A:ASP:H	1:144:A:ASN:HA	2	0.32	0.04	0.32

¹Number of violated models, ²Standard deviation

9.5 All violated distance restraints [i](#)

9.5.1 Histogram : Distribution of distance violations [i](#)

The following histogram shows the distribution of the absolute value of the violation for all violated restraints in the ensemble.



9.5.2 Table : All distance violations [i](#)

The following table lists the absolute value of the violation for each restraint in the ensemble sorted by its value. The Key (restraint list ID, restraint ID) is the unique identifier for a given restraint. Rows with same key represent combinatorial or ambiguous restraints and are counted as a single restraint.

Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,6)	1:181:A:ALA:H	1:157:A:ALA:HA	5	0.84
(1,4)	1:150:A:TRP:H	1:92:A:LEU:HG	3	0.64
(1,29)	1:122:A:THR:H	1:145:A:ALA:HA	4	0.5
(1,33)	1:133:A:ASP:H	1:144:A:ASN:HA	4	0.36
(1,33)	1:133:A:ASP:H	1:144:A:ASN:HA	9	0.27
(1,29)	1:122:A:THR:H	1:145:A:ALA:HA	9	0.23
(1,6)	1:181:A:ALA:H	1:157:A:ALA:HA	9	0.13
(1,25)	1:118:A:VAL:H	1:111:A:SER:HB2	10	0.11
(1,25)	1:118:A:VAL:H	1:111:A:SER:HB3	10	0.11

10 Dihedral-angle violation analysis ⓘ

No dihedral-angle restraints found